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NEWS	4	NOV		CHEMSAFE now available on STN Easy
	-			
NEWS	5	NOV		Two new SET commands increase convenience of STN searching
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			-	Classification Data
NEWS	11	FEB	02	Simultaneous left and right truncation (SLART) added for CERAB, COMPUAB, ELCOM, and SOLIDSTATE
NEWS	12	FEB	02	GENBANK enhanced with SET PLURALS and SET SPELLING
NEWS	13	FEB	06	Patent sequence location (PSL) data added to USGENE
NEWS				COMPENDEX reloaded and enhanced
NEWS		FEB		WTEXTILES reloaded and enhanced
NEWS	16	FEB	19	New patent-examiner citations in 300,000 CA/CAplus patent records provide insights into related prior art
NEWS	17	FEB	19	Increase the precision of your patent queries use terms from the IPC Thesaurus, Version 2009.01
NEWS	18	FEB	23	Several formats for image display and print options discontinued in USPATFULL and USPAT2
NEWS	19	FEB	23	MEDLINE now offers more precise author group fields and 2009 MeSH terms
NEWS	20	FEB	23	TOXCENTER updates mirror those of MEDLINE - more precise author group fields and 2009 MeSH terms
NEWS	21	FEB	23	Three million new patent records blast AEROSPACE into STN patent clusters
NEWS	22	FEB	25	USGENE enhanced with patent family and legal status display data from INPADOCDB
NEWS	23	MAR	06	INPADOCDB and INPAFAMDB enhanced with new display formats
NEWS	24	MAR	11	EPFULL backfile enhanced with additional full-text
	0.5			applications and grants
NEWS		MAR		ESBIOBASE reloaded and enhanced
NEWS	26	MAR	20	CAS databases on STN enhanced with new super role

for nanomaterial substances

NEWS 27 MAR 23 CA/Caplus enhanced with more than 250,000 patent equivalents from China

NEWS 28 MAR 30 IMSPATENTS reloaded and enhanced

NEWS 29 APR 03 CAS coverage of exemplified prophetic substances enhanced

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Uploading C:\Program Files\STNEXP\Queries\10551737 R5 aryl R8 and R9 ring.str

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7 8 12 13 14 17 19 20
ring nodes :
1 2 3 4 5 6 10 11
chain bonds :
4-17 5-7 7-8 8-20 10-20 11-12 12-13 12-14 17-19
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 10-11
exact/norm bonds :
4-17 5-7 7-8 10-11 10-20 12-13 12-14 17-19
exact bonds :
8-20 11-12
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6
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G1:C.O.S

G2:0,S

G3:Cb,Cy,Hy

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 10:CLASS 12:CLASS 13:CLASS 14:CLASS 17:CLASS 19:CLASS 20:CLASS

## L1 STRUCTURE UPLOADED

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Structure attributes must be viewed using STN Express query preparation.

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FULL ESTIMATED COST 0.48 0.70

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=> s L1 SSS full REGISTRY INITIATED

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FULL SEARCH INITIATED 07:38:38 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 184852 TO ITERATE

100.0% PROCESSED 184852 ITERATIONS SEARCH TIME: 00.00.11

65 ANSWERS

L2 65 SEA SSS FUL L1

L3 4 L2

=> d ibib abs hitstr 1-YOU HAVE REQUESTED DATA FROM 4 ANSWERS - CONTINUE? Y/(N):y

L3 ANSWER 1 OF 4 CAPLUS COPYRIGHT 2009 ACS on STN ACCESSION NUMBER: 2006:1093266 CAPLUS Full-text

DOCUMENT NUMBER: 145:432223

TITLE: Method of treating schizophrenia prodrome

INVENTOR(S): Woods, Scott W.
PATENT ASSIGNEE(S): Yale University, USA
SOURCE: PCT Int. Appl., 64pp.
CODEN: PIXXD.

DOCUMENT TYPE: Patent

LANGUAGE: English FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

			KIND DATE				APPLICATION NO.											
WO	2006110724				A2		20061019											
WO	WO 2006110724 W: AE, AG, AL,																	
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		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KΕ,	KG,	KM,	KN,	KP,	KR,	
		KZ,	LC,	LK,	LR,	LS,	LT,	LU,	LV,	LY,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	
		MZ,	NA,	NG,	NI,	NO,	NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	
		SG,	SK,	SL,	SM,	SY,	TJ,	TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UZ,	VC,	
		VN,	YU,	ZA,	ZM,	ZW												
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PRIORIT	Y APP	LN.	INFO	.:						US 2005-670600P								
										WO 2	006-	US13	444		W 2	0060	411	
OTHER S	OTHER SOURCE(S):						145:	43222	23									

- The present invention relates to a method of treating schizophrenia prodrome AB in human subjects using a NMDA glycine site agonist, a glycine transporter-1 inhibitor or mixts. thereof, optionally in combination with a pharmaceutically acceptable additive, carrier or excipient.
- 791642-83-6

RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

(method of treating schizophrenia prodrome with NMDA glycine agonist and glycine transporter-1 inhibitor)

- 791642-83-6 CAPLUS RN
- CN L-Proline, 1-[2-[2-[[4-(1,1-dimethylethyl)phenyl]thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS REFERENCE COUNT: RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

ANSWER 2 OF 4 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2006:625349 CAPLUS Full-text

DOCUMENT NUMBER: 145:224321

TITLE: The synthesis and SAR of 2-arylsulfanylphenyl-1-oxyalkylamino acids as GlyT-1

inhibitors Smith, Garrick; Mikkelsen, Gitte; Eskildsen, Jorgen;

AUTHOR(S): Bundgaard, Christoffer

CORPORATE SOURCE: Medicinal Chemistry Research, H. Lundbeck A/S, Valby,

DK 2500, Den.

SOURCE: Bioorganic & Medicinal Chemistry Letters (2006),

16(15), 3981-3984

CODEN: BMCLE8: ISSN: 0960-894X

Elsevier B.V.

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 145:224321

GI

PUBLISHER:

- AB Elevation of glycine levels by inhibition of the glycine transporter-1 [GlyT-1] and activation of the NMDA receptor is a potential strategy for the treatment of schizophrenia. A novel series of 2-arylsulfanylphenyl-1-oxyalkyl amino acids have been identified. The most prominent member of this series [I) is a potent GlyT-1 inhibitor (ICSD = 59 nM). In vitro and in vivo assessment of CNS exposure indicates this compound is a likely substrate for active efflux transporters.
- IT 791644-20-7P 791644-21-0P
  RL: BSU (Biological study, unclassified); PAC (Pharmacological activity);
  SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
  (synthesis and SAR of arylsulfanylphenyloxyalkylamino acids as GlyI-1 inhibitors)

Absolute stereochemistry.

- RN 791644-21-8 CAPLUS
- CN L-Proline, 1-[2-[[3-[(3-fluorophenyl)thio]-4'-methoxy[1,1'-biphenyl]-4yl]oxylethyl]- (CA INDEX NAME)

905815-64-7P 905815-65-8P 905815-66-9P

905815-67-0P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis and SAR of arylsulfanylphenyloxyalkylamino acids as GlyT-1 inhibitors)

RN 791642-87-0 CAPLUS

CN L-Proline, 1-[2-[2-[(3-chlorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-17-2 CAPLUS

N L-Proline, 1-[2-[3-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-18-3 CAPLUS

CN L-Proline, 1-[2-[5-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

RN 794510-03-5 CAPLUS

CN L-Proline, 1-[2-[4-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 905815-62-5 CAPLUS

CN D-Proline, 1-[2-[2-[(3-chlorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 905815-63-6 CAPLUS

CN L-Proline, 1-[2-[2-chloro-6-[(3-fluoropheny1)thio]phenoxy]ethyl]- (CA INDEX NAME)

RN 905815-64-7 CAPLUS

Absolute stereochemistry.

RN 905815-65-8 CAPLUS

CN L-Proline, 1-[2-[2-[(3-fluorophenyl)thio]-4-(3-thienyl)phenoxy]ethyl] (CA INDEX NAME)

Absolute stereochemistry.

RN 905815-66-9 CAPLUS

CN L-Proline, 1-[2-[[3-fluorophenyl)thio]-3'-methoxy[1,1'-biphenyl]-4yl]oxy]ethyl]- (CA INDEX NAME)

- RN 905815-67-0 CAPLUS
- CN L-Proline, 1-[2-[[4'-chloro-3-[(3-fluorophenyl)thio][1,1'-biphenyl]-4yl]oxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

- IT 791642-79-0P 791644-01-4P 905816-02-6P
  - 905816-03-7P 905816-06-0P 905816-07-1P
  - 905816-08-2P 905816-09-3P
  - RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
  - (synthesis and SAR of arylsulfanylphenyloxyalkylamino acids as GlyT-1 inhibitors)
- RN 791642-79-0 CAPLUS
- CN L-Proline, 1-[2-[4-bromo-2-[(3-fluorophenyl)thio]phenoxy]ethyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

RN 791644-01-4 CAPLUS

CN L-Proline, 1-[2-[5-bromo-2-[(3-fluoropheny1)thio]phenoxy]ethyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

Absolute stereochemistry.

RN 905816-02-6 CAPLUS

CN D-Proline, 1-[2-[2-[(3-chlorophenyl)thio]phenoxy]ethyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

Absolute stereochemistry.

RN 905816-03-7 CAPLUS

CN L-Proline, 1-[2-[2-[(3-chlorophenyl)thio]phenoxy]ethyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

RN 905816-06-0 CAPLUS

CN L-Proline, 1-[2-[5-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

Absolute stereochemistry.

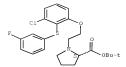
RN 905816-07-1 CAPLUS

CN L-Proline, 1-[2-[2-chloro-6-[(3-fluorophenyl)thio]phenoxy]ethyl]-,
1,1-dimethylethyl ester (CA INDEX NAME)

Absolute stereochemistry.

RN 905816-08-2 CAPLUS

CN L-Proline, 1-[2-[3-chloro-2-[(3-fluoropheny1)thio]phenoxy]ethyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)



905816-09-3 CAPLUS

L-Proline, 1-[2-[4-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

Absolute stereochemistry.

REFERENCE COUNT: 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L3 ANSWER 3 OF 4 CAPLUS COPYRIGHT 2009 ACS on STN ACCESSION NUMBER: 2004:965214 CAPLUS Full-text

DOCUMENT NUMBER: 141:411217

TITLE: A preparation of oxyphenyl and sulfanylphenyl derivatives of amino acids, useful as glycine

transporter inhibitors

INVENTOR(S): Smith, Garrick Paul; Mikkelsen, Gitte; Andersen, Kim;

Greve, Daniel Rodriguez; Eskildsen, Joergen

PATENT ASSIGNEE(S): H. Lundbeck A/S, Den. SOURCE: PCT Int. Appl., 87 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.						D	DATE			APPL	ICAT	ION :	DATE					
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WO 2004096761							20041111			WO 2	004-	20040427						
	W:	ΑE,	AG,	AL,	AM,	AT,	AU,	AZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,	
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FΙ,	GB,	GD,	
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	ΚZ,	LC,	
		LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NA,	NI,	
		NO.	NZ.	OM.	PG.	PH.	PL.	PT.	RO.	RU.	SC.	SD.	SE.	SG.	SK.	SL.	SY.	

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             EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE,
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                                                                    20030430
                                             US 2003-466755P
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                                             WO 2004-DK290
                                                                    20040427
OTHER SOURCE(S):
                         MARPAT 141:411217
```

AB The invention relates to a preparation of aromatic oxyphenyl and aromatic sulfanylphenyl derivs. of formula I [wherein: X is O, S, or CH2, etc.; Y is O or S; R1, R2, R3, and R4 are independently selected from H, halogen, CN, NO2, or alk(en/yn)yl, etc.; R5 is (un)substituted aryl or monocyclic heteroaryl; R6 is H, alk(en/yn)yl, cycloalk(en)yl, or alk(en/yn)ylsulfanyl, etc.; R7 and R8 are independently selected from H, alk(en/yn)yl, or cycloalk(en)yl; R9 and R11 are independently selected from H, alk(en/yn)yl, hydroxyalk(en/yn)yl, or alk(en/yn)ylsulfanyl, etc.; R10 is H, alk(en/yn)yl, aryl, or arylalk(en/yn)yl, etc.; R6 and R8 together with the nitrogen may form 3-7 membered heterocyclic ring], useful as glycine transporter inhibitors (IC50 < 10000 nM). The compds. of formula I are useful for the treatment of diseases such as schizophrenia, including both the pos. and the neg. symptoms of schizophrenia. For instance, pyrrolidinecarboxylic acid derivative II was prepared via etherification of 2-(3-fluorophenylsulfanyl)phenol by (hydroxyethyl)pyrrolidinecarboxylate derivative III.

IT 791642-79-9F, (S)-1-[2-[4-Bromo-2-(3-fluorophenylsulfanyl)phenoxylethyl]pyrrolidine-2-carboxylic acid

tert-butyl ester RI: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(intermediate; preparation of oxyphenyl and sulfanylphenyl derivs. of amino acids, useful as glycine transporter inhibitors)

RN 791642-79-0 CAPLUS

CN L-Proline, 1-[2-[4-bromo-2-[(3-fluoropheny1)thio]phenoxy]ethyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

Absolute stereochemistry.

```
791642-81-4P, (S)-1-[2-[2-(4-
Fluorophenylsulfanyl)phenoxy]ethyl]pyrrolidine-2-carboxylic acid
791642-83-6P, (S)-1-[2-[2-(4-tert-
Butylphenylsulfanyl)phenoxy]ethyl]pyrrolidine-2-carboxylic acid
791642-84-7P, (S)-1-[2-[2-(4-
Trifluoromethylphenylsulfanyl)phenoxy]ethyl]pyrrolidine-2-carboxylic acid
791642-85-8P, (S)-1-[2-[2-(3-
Fluorophenylsulfanyl)phenoxy]ethyl]pyrrolidine-2-carboxylic acid
791642-86-9P, (S)-1-[2-[2-(4-Chlorophenylsulfanyl)-phenoxy]-
ethyl]pyrrolidine-2-carboxylic acid 791642-87-0P,
(S)-1-[2-[2-(3-Chlorophenylsulfanyl)phenoxy]ethyl]pyrrolidine-2-carboxylic
acid 791642-88-1P, (S)-1-[2-[2-(3,4-
Dichlorophenylsulfanyl)phenoxy]ethyl]pyrrolidine-2-carboxylic acid
791642-90-5P, (S)-1-[2-[2-(3-Chloro-4-
fluorophenylsulfanyl)phenoxy]ethyl]pyrrolidine-2-carboxylic acid
791642-91-6P, (S)-1-[2-[2-(3-
Chlorophenoxy)phenoxy]ethyl]pyrrolidine-2-carboxylic acid
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791644-28-5P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
(Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
(Uses)
```

(preparation of oxyphenyl and sulfanylphenyl derivs. of amino acids, useful

as glycine transporter inhibitors)

RN 791642-81-4 CAPLUS

CN L-Proline, 1-[2-[2-[(4-fluorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-83-6 CAPLUS

CN L-Proline, 1-[2-[2-[[4-(1,1-dimethylethyl)phenyl]thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-84-7 CAPLUS

CN L-Proline, 1-[2-[2-[[4-(trifluoromethyl)phenyl]thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-85-8 CAPLUS

CN L-Proline, 1-[2-[3-fluorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-86-9 CAPLUS

CN L-Proline, 1-[2-[2-[(4-chlorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-87-0 CAPLUS

CN L-Proline, 1-[2-[2-[(3-chlorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-88-1 CAPLUS

CN L-Proline, 1-[2-[2-[(3,4-dichlorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

RN 791642-90-5 CAPLUS

CN L-Proline, 1-[2-[2-[(3-chloro-4-fluorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-91-6 CAPLUS

CN L-Proline, 1-[2-[2-(3-chlorophenoxy)phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-92-7 CAPLUS

CN L-Proline, 1-[2-[2-(4-chlorophenoxy)phenoxy]ethyl]- (CA INDEX NAME)

RN 791642-93-8 CAPLUS

CN L-Proline, 1-[2-[2-(4-methoxyphenoxy)phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-94-9 CAPLUS

CN L-Proline, 1-[2-[2-(3,4-difluorophenoxy)phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-95-0 CAPLUS

CN L-Proline, 1-[2-[2-(4-chlorophenoxy)phenoxy]propyl]- (CA INDEX NAME)

RN 791642-97-2 CAPLUS

CN L-Proline, 1-[2-[2-(3,4-difluorophenoxy)phenoxy]propyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-98-3 CAPLUS

CN L-Proline, 1-[2-[2-(3-fluorophenoxy)phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791642-99-4 CAPLUS

CN L-Proline, 1-[2-[2-(3-fluorophenoxy)phenoxy]propyl]- (CA INDEX NAME)

RN 791643-00-0 CAPLUS

CN L-Proline, 1-[2-[2-[(3-fluorophenyl)thio]phenoxy]propyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791643-01-1 CAPLUS

CN L-Proline, 1-[2-[2-[(3-chlorophenyl)thio]phenoxy]propyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791643-85-1 CAPLUS

CN L-Proline, 1-[3-[2-[(3-fluorophenyl)thio]phenyl]propyl]-, hydrochloride (9CI) (CA INDEX NAME)

● HCl

RN 791643-88-4 CAPLUS

CN L-Proline, 1-[2-[4-chloro-2-[(3-fluoropheny1)thio]phenoxy]ethyl]-, hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HCl

RN 791643-90-8 CAPLUS

CN L-Proline, 1-[2-[3-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]-,
hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

RN 791643-91-9 CAPLUS

CN L-Proline, 1-[2-[5-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]-, hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

- RN 791643-92-0 CAPLUS
- CN L-Proline, 1-[2-[4-cyano-2-[(3-fluoropheny1)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

- RN 791643-94-2 CAPLUS
- CN L-Proline, 1-[2-[5-chloro-2-(phenylthio)phenoxy]ethyl]-, hydrochloride (9CI) (CA INDEX NAME)

RN 791643-95-3 CAPLUS
CN L-Proline, 1-[2-[[3-[(3-fluorophenyl)thio][1,1'-biphenyl]-4-yl]oxy]ethyl]-

. hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HC1

RN 791643-97-5 CAPLUS

CN L-Proline, 1-[2-[[3-[(3-fluorophenyl)thio]-4'-methoxy[1,1'-biphenyl]-4yl]oxy]ethyl]-, hydrochloride (9CI) (CA INDEX NAME)

RN 791643-99-7 CAPLUS

CN L-Proline, 1-[2-[[4'-cyano-3-[(3-fluorophenyl)thio][1,1'-biphenyl]-4-yl]oxy]ethyl]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-00-3 CAPLUS

CN L-Proline, 1-[2-[[4'-cyano-4-[(3-fluorophenyl)thio][1,1'-biphenyl]-3yl]oxy]ethyl]-, monohydrochloride (9CI) (CA INDEX NAME)

RN 791644-02-5 CAPLUS

CN L-Proline, 1-[2-[2-[(3-fluoropheny1)thio]-5-(3-thieny1)phenoxy]ethyl]-, hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-04-7 CAPLUS

CN L-Proline, 1-[2-[2-[(3-fluoropheny1)thio]-4-(5-pyrimidiny1)phenoxy]ethy1]-, monohydrochloride (9CI) (CA INDEX NAME)

● HCl

RN 791644-06-9 CAPLUS
CN L-Proline, 1-[2-[[3-[(3-fluorophenyl)thio]-3'-(methylsulfonyl)[1,1'-biphenyl]-4-ylloxy[sthyl]-, hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-08-1 CAPLUS

CN L-Proline, 1-[2-[2-[(3-fluorophenyl)thio]-4-(4-morpholinyl)phenoxy]ethyl]-, monohydrochloride (9CI) (CA INDEX NAME)

HC1

RN 791644-09-2 CAPLUS

CN L-Proline, 1-[2-[2-[(3-fluorophenyl)thio]-4-(1-piperidinyl)phenoxy]ethyl]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-15-0 CAPLUS

HC1

CN L-Proline, 1-[3-[2-[(3-fluorophenyl)thio]phenyl]propyl]- (CA INDEX NAME)

RN 791644-17-2 CAPLUS

CN L-Proline, 1-[2-[3-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-18-3 CAPLUS

CN L-Proline, 1-[2-[5-chloro-2-[(3-fluorophenyl)thio]phenoxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-19-4 CAPLUS

CN L-Proline, 1-[2-[5-chloro-2-(phenylthio)phenoxy]ethyl]- (CA INDEX NAME)

RN 791644-20-7 CAPLUS

CN L-Proline, 1-[2-[[3-[(3-fluorophenyl)thio][1,1'-biphenyl]-4-yl]oxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

- RN 791644-21-8 CAPLUS
- CN L-Proline, 1-[2-[[3-[(3-fluorophenyl)thio]-4'-methoxy[1,1'-biphenyl]-4yl]oxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

- RN 791644-22-9 CAPLUS
- CN L-Proline, 1-[2-[[4'-cyano-3-[(3-fluorophenyl)thio][1,1'-biphenyl]-4yl]oxy]ethyl]- (CA INDEX NAME)

RN 791644-23-0 CAPLUS

CN L-Proline, 1-[2-[[4'-cyano-4-[(3-fluorophenyl)thio][1,1'-biphenyl]-3yl]oxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-24-1 CAPLUS

CN L-Proline, 1-[2-[2-[(3-fluorophenyl)thio]-5-(3-thienyl)phenoxy]ethyl]-(CA INDEX NAME)

Absolute stereochemistry.

RN 791644-25-2 CAPLUS

RN 791644-26-3 CAPLUS

CN L-Proline, 1-[2-[[3-[(3-fluorophenyl)thio]-3'-(methylsulfonyl)[1,1'-biphenyl]-4-yl]oxy]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 791644-27-4 CAPLUS

Absolute stereochemistry.

- IT 791643-98-6 791644-01-4 791644-07-0
  - RL: RCT (Reactant); RACT (Reactant or reagent)
    (reactant; preparation of oxyphenyl and sulfanylphenyl derivs. of amino
- acids, useful as glycine transporter inhibitors)
  RN 791643-98-6 CAPLUS
- CN L-Proline, 1-[2-[4-bromo-2-[(3-fluorophenyl)thio]phenoxy]ethyl]-, hydrochloride (901) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

- RN 791644-01-4 CAPLUS
- CN L-Proline, 1-[2-[5-bromo-2-[(3-fluorophenyl)thio]phenoxy]ethyl]-,
  1,1-dimethylethyl ester (CA INDEX NAME)

RN 791644-07-0 CAPLUS

L-Proline, 1-[2-[4-bromo-2-[(3-fluorophenyl)thio]phenoxy]ethyl]-, butyl CN ester (CA INDEX NAME)

Absolute stereochemistry.

REFERENCE COUNT: THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L3 ANSWER 4 OF 4 CAPLUS COPYRIGHT 2009 ACS on STN ACCESSION NUMBER: 2000:666715 CAPLUS Full-text

DOCUMENT NUMBER: 133:252449

TITLE: Quinazolines and other bicyclic heterocycles,

pharmaceutical compositions containing these compounds as tyrosine kinase inhibitors, and processes for

preparing them

INVENTOR(S): Himmelsbach, Frank; Langkopf, Elke; Blech, Stefan;

Jung, Birgit; Metz, Thomas; Solca, Flavio PATENT ASSIGNEE(S): Boehringer Ingelheim Pharma K.-G., Germany

PCT Int. Appl., 153 pp.

SOURCE: CODEN: PIXXD2

DOCUMENT TYPE: Patent English

LANGUAGE: FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.						KIN	D	DATE			APPL	CAT	DATE						
	WO 2000055141					A1 20000921					WO 2	000-	20000314						
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			CZ,	DE,	DK,	DM,	EE,	ES,	FΙ,	GB,	GD,	GE,	GH,	GM,	HR,	HU,	ID,	IL,	
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	3754				B2		2006	0315										
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	1150				C			0519		2000-					0000			
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	2689				В			1221	TW	2000-		20000426						
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OTHER SOURCE(S): MARPAT 133:252449

The invention relates to bicyclic heterocyclic compds. I [R1 = H, alkyl; R2 = AB (un) substituted Ph, CH2Ph, or CH(Me)Ph; R3, R4 = H, F, C1, OMe, or Me optionally substituted by OMe, NMe2, NEt2, pyrrolidino, piperidino, or morpholino; X = N or C(CN); A = O, NH, (un)substituted alkylene, O-alkylene, NH-alkylene, O-cycloalkylene, etc.; B = (un) substituted amine-containing sidechain, piperazino, alkyleneimino, morpholino, etc.; or AB = H, F, Cl, alkoxy, amino, etc.; C = groups similar to A; D = groups similar to B; with a variety of provisos | and their tautomers, stereoisomers, and salts, and particularly their physiol, acceptable salts with inorg, or organic acids or bases. The compds, have valuable pharmacol, properties, particularly an inhibitory effect on signal transduction mediated by tyrosine kinases, and are useful in treating diseases, particularly tumor diseases, and diseases of the lung and airways. Over 20 compds. were prepared, and over 200 are listed. For instance, alkylation of 4-(3-chloro-4-fluorophenylamino)-6-[3-(1piperazinyl)propyloxyl-7- methoxyquinazoline (preparation given) by Me bromoacetate gave 51% title compound II. The latter compound inhibited EGFdependent proliferation of F/L-HERC cells in vitro, with an IC50 of 46 nM. IT 295330-27-7P, (R)-4-[(3-Chloro-4-fluorophenyl)amino]-6-[2-[2-

(methoxycarbonyl)pyrrolidin-1-yl]ethoxy]-7-cyclopentyloxyquinazoline
RL: BAC (Biological activity or effector, except adverse); BSU (Biological
study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
BIOL (Biological study); PREP (Preparation); USES (Uses)

(drug candidate; preparation of quinazoline derivs. and other bicyclic heterocycles as tyrosine kinase inhibitors)

RN 295330-27-7 CAPLUS

CN D-Proline, 1-[2-[[4-[(3-chloro-4-fluorophenyl)amino]-7-(cyclopentyloxy)-6quinazolinyl]oxy]ethyl]-, methyl ester (CA INDEX NAME)

REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD, ALL CITATIONS AVAILABLE IN THE RE FORMAT

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